

Sum rules of single-particle spectral functions in hot asymmetric nuclear matter

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The neutron and proton single-particle spectral functions in asymmetric nuclear matter fulfill energy weighted sum rules. The validity of these sum rules within the self-consistent Green's function approach is investigated. The various contributions to these sum rules and their convergence as a function of energy provide information about correlations induced by the realistic interaction between the nucleons. These features are studied as a function of the asymmetry of nuclear matter.

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I. INTRODUCTION

The microscopic description of the single-particle properties in nuclear matter must deal with the treatment of nucleon-nucleon (NN) correlations [1, 2]. In fact, the strong short-range and tensor components which are required in realistic NN interactions to fit the NN scattering data, lead to corresponding correlations in the nuclear wave function. These NN correlations are important to describe bulk properties of dense matter. They also modify the spectral distribution of the single-particle strength in a significant way. Recent calculations have demonstrated without ambiguity that NN correlations produce a partial occupation of the single-particle states, which are completely occupied in the mean field approach, and a wide distribution of the single-particle strength in energy. These two features have found experimental grounds in the analysis of $(e, e'p)$ reactions [3, 4].

Historically, various tools have been employed to account for correlations in the nuclear many-body wave function. These include the traditional Brueckner hole-line expansion [5] and variational approaches using correlated basis functions [6]. Attempts have also been made to employ the technique of a self-consistent evaluation of Green's functions (SCGF) [1, 2, 7, 8] to the solution of the nuclear many-body problem. This method offers various advantages: (i) the single-particle Greens function contains detailed information about the spectral function, i.e., the distribution of single-particle strength as a function of missing energy and momentum; (ii) the method can be extended to finite temperatures, a feature which is of interest for the study of the nuclear properties in astrophysical environments; (iii) the Brueckner Hartree-Fock (BHF) approximation, the approximation to the hole-line expansion which is commonly used, can be considered as a specific prescription within this scheme.

Enormous progress in the SCGF applications to nuclear matter has been achieved in the last years, both at zero [9] and finite temperature [10, 11, 12, 13]. The efforts at $T = 0$ are mainly oriented to provide the appropriate theoretical support for the interpretation of the $(e, e'p)$ experiments, while research at finite T is essentially focused on the description of the nuclear medium in astrophysical environments or on the interpretation of heavy ion collisions dynamics. In all cases, a key quantity is the single-particle spectral function which measures the possibility to add or remove a particle with a given momentum at a specific energy. A useful way to study the properties of these single-particle spectral functions is by means of the energy weighted sum rules, which are well established in the literature and which have been numerically analyzed for symmetric nuclear matter at zero [14] and finite temperature [15].

Recently, the single-particle spectral functions in hot asymmetric nuclear matter have been calculated within the SCGF framework [13]. In this computation, the bulk properties of asymmetric dense matter were computed, namely the energy per particle, the symmetry energy and the chemical potentials. In addition, and since the SCGF approach gives access to the correlated momentum distributions of neutrons and protons, the influence of the asymmetry on the depletion of the momentum states could be studied. In neutron-rich matter the proton states with momenta below the Fermi momentum are more strongly depleted than in the symmetric system at the same density. In this case the occupation of neutron states within their Fermi sphere are enhanced compared to the symmetric case. This

indicates that the proton-neutron interaction is a stronger source of correlation as compared to proton-proton and neutron-neutron interaction.

In this paper, we want to use the energy weighted sum rules to investigate the dependence of the single-particle spectral functions on the asymmetry more in detail. These sum rules may help to explore the isospin dependence of the short-range correlations in asymmetric matter. Besides, the analysis of the energy weighted sum rules provides valuable tests on the numerical accuracy of our calculations. All the computations discussed in this paper are performed in the framework of SCGF employing a fully self-consistent ladder approximation in which the complete spectral function has been used to describe the intermediate states in the Galitskii-Feynman equation [12, 13].

After this introduction, we outline the derivation of the energy weighted sum rules in Section 2. The results obtained for hot asymmetric nuclear using the charge dependent Bonn potential CDBONN [16] are presented in Section 3, where a short summary of the main conclusions is given as well.

II. SUM RULES FOR ASYMMETRIC NUCLEAR MATTER

For a given Hamiltonian H , the single-particle Green's function for a system at finite temperature can be defined in a grand-canonical formulation:

$$ig_\tau(\mathbf{k}t; \mathbf{k}'t') = \text{Tr} \left\{ \rho \mathbf{T} \left[a_{\mathbf{k}\tau}(t) a_{\mathbf{k}'\tau}^\dagger(t') \right] \right\} , \quad (1)$$

where the subindex $\tau = n(p)$ refers either to neutrons or protons and \mathbf{T} is the time ordering operator that acts on a product of Heisenberg field operators $a_{\mathbf{k}\tau}(t) = e^{itH} a_{\mathbf{k}\tau} e^{-itH}$, in such a way that the field operator with the largest time argument is put to the left. The trace is to be taken over all energy eigenstates and all particle number eigenstates of the many-body system, weighted by the statistical operator,

$$\rho = \frac{1}{Z} e^{-\beta(H - \mu_n N_n - \mu_p N_p)} . \quad (2)$$

β denotes the inverse of the temperature and μ_n and μ_p are the neutron and proton chemical potentials associated to a given average number of neutrons and protons. N_n and N_p are the operators that count the total number of neutrons and protons in the system,

$$N_\tau = \sum_{\mathbf{k}} a_{\mathbf{k}\tau}^\dagger(t) a_{\mathbf{k}\tau}(t) . \quad (3)$$

The normalization factor in Eq. (2) is the grand partition function of statistical mechanics,

$$Z = \text{Tr} \, e^{-\beta(H - \mu_n N_n - \mu_p N_p)} . \quad (4)$$

For a homogeneous system, the Green's function is diagonal in momentum space and depends only on the absolute value of \mathbf{k} and on the difference $\xi = t' - t$. Starting from the definition of the Green's function, we first focus on the case $\xi > 0$. In order to recover the expression in the grand canonical ensemble average of the occupation number $n_\tau(k)$ for $\xi = 0^+$, the following definition of the correlation functions $g_\tau^<$ includes an additional factor $-i$ with respect to the definitions of the Green's function g_ν in Eq. (1),

$$g_\nu^<(k, \xi) = \text{Tr} \left\{ \rho e^{i\xi H} a_{\mathbf{k}\nu}^\dagger e^{-i\xi H} a_{\mathbf{k}\nu} \right\} . \quad (5)$$

$g_\nu^<(k, \xi)$ can be expressed as a Fourier integral over the frequencies,

$$g_\nu^<(k, \xi) = \int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} e^{-i\omega\xi} A_\nu^<(k, \omega) \quad (6)$$

where $A_\nu^<(k, \omega)$ is defined by

$$A_\nu^<(k, \omega) = 2\pi \sum_{rs} \frac{e^{-\beta(E_r - \mu_p N_{r(p)} - \mu_n N_{r(n)})}}{Z} |\langle \Psi_s | a_{\mathbf{k}\nu} | \Psi_r \rangle|^2 \delta(\omega - (E_r - E_s)) . \quad (7)$$

A similar analysis can be conducted for $\xi < 0$, yielding a function

$$A_\nu^>(k, \omega) = e^{\beta(\omega - \mu_\nu)} A_\nu^<(k, \omega) . \quad (8)$$

Finally, the single-particle spectral function is defined as the sum of the two positive functions, $A_\nu^<$ and $A_\nu^>$,

$$A_\nu(k, \omega) = A_\nu^<(k, \omega) + A_\nu^>(k, \omega) . \quad (9)$$

In our computations, we obtain the full spectral function $A_\nu(k, \omega)$ and we make use of the relations:

$$A_\nu^<(k, \omega) = f(\omega) A_\nu(k, \omega) , \quad A_\nu^>(k, \omega) = (1 - f(\omega)) A_\nu(k, \omega) \quad (10)$$

to obtain $A_\nu^<$ and $A_\nu^>$, where $f(\omega)$ is the Fermi-Dirac distribution function. The integration of the function $A_\nu^<$ leads to the occupation probability

$$n_\nu(k) = \int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} A_\nu^<(k, \omega) , \quad (11)$$

for the state with isospin ν and momentum k at the inverse temperature β .

The functions $A_\nu^<$ and $A_\nu^>$ can be compared with the hole and particle part of the spectral function at zero temperature. In particular, the limit of $A_\nu^<$ at zero temperature reads

$$A_\nu^h = 2\pi \sum_r |\langle \Psi_r^{A-1} | a_{k\nu} | \Psi_0^A \rangle|^2 \delta(\omega - (E_0^A - E_r^{A-1})), \quad (12)$$

where $|\Psi_0^A\rangle$ is the ground state with N_n neutrons and N_p protons (such that $A = N_p + N_n$) and $|\Psi_r^{A-1}\rangle$ labels the excited energy state of a system with a neutron or a proton less (depending on which type of annihilation operator $a_{k\nu}$ has been applied to the ground state). By its own definition, it is clear that the lowest possible energy of the final state is the ground state energy of the $A - 1$ particle system, so that there is an upper limit for the hole spectral function $\omega_\nu = E_0^A - E_0^{A-1} = \mu_\nu$, both for neutrons and protons. In a similar way one can define the particle part of the spectral function and find a lower bound for the excitation energy of the $A + 1$ particle system, measured with respect to the ground state of the A particle system. At zero temperature, the existence of these lower and upper bounds causes a complete separation in energy between the particle and the hole part of the spectral function.

In the T matrix approximation to the self-energy reported in Ref.[13], one can determine the single-particle Green's function as the solution of Dyson's equation for any complex value of the frequency variable z

$$g_\nu(k, z) = \frac{1}{z - \frac{k^2}{2m} - \Sigma_\nu(k, z)}. \quad (13)$$

The analytical properties of the finite temperature Green's function allows one to derive the corresponding Lehmann representation, which for slightly complex values of the frequency can be written as

$$g_\nu(k, \omega + i\eta) = \int_{-\infty}^{+\infty} \frac{d\omega'}{2\pi} \frac{A(k, \omega')}{\omega - \omega' + i\eta}. \quad (14)$$

The sum rules for the spectral functions can be derived from the asymptotic behavior at large ω by expanding the real part of both previous expressions for the Green's function, Eqs. (13) and (14), in powers of $1/\omega$. This yields

$$\text{Re } g_\nu(k, \omega) = \frac{1}{\omega} \left\{ 1 + \frac{1}{\omega} \left[\frac{k^2}{2m} + \lim_{\omega \rightarrow \infty} \text{Re} \Sigma_\nu(k, \omega) \right] + \dots \right\} \quad (15)$$

and

$$\text{Re } g_\nu(k, \omega) = \frac{1}{\omega} \left\{ \int_{-\infty}^{+\infty} d\omega' A_\nu(k, \omega') + \frac{1}{\omega} \int_{-\infty}^{+\infty} d\omega' \omega' A_\nu(k, \omega') + \dots \right\}. \quad (16)$$

By comparing the first two expansion coefficients, one finds the m_0

$$\int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} A_\nu(k, \omega) = 1, \quad (17)$$

and the m_1 sum rules

$$\int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} \omega A_\nu(k, \omega) = \frac{k^2}{2m} + \lim_{\omega \rightarrow \infty} \text{Re} \Sigma_\nu(k, \omega). \quad (18)$$

It is worth to mention that in our scheme the self-energy is derived in the T matrix approximation and so its real part is computed from the imaginary part using the following dispersion relation:

$$\text{Re} \Sigma_\nu(k, \omega) = \Sigma_\nu^\infty(k) - \frac{\mathcal{P}}{\pi} \int_{-\infty}^{+\infty} d\lambda \frac{\text{Im} \Sigma_\nu(k, \lambda + i\eta)}{\omega - \lambda}. \quad (19)$$

In the derivation of the previous equation, the spectral decomposition of the Green's function is already used, so it is a property of the T matrix approach that it fulfils the sum rule. Nevertheless, the sum rules still provide a useful consistency check for the numerics. The first term on the right hand side of Eq. (19) is the energy independent part of the self-energy,

$$\Sigma_\nu^\infty(k) = \sum_\tau \int \frac{d^3 k'}{(2\pi)^3} \langle \mathbf{k} \nu \mathbf{k}' \tau | V | \mathbf{k} \nu \mathbf{k}' \tau \rangle_A n_\tau(\mathbf{k}') \quad (20)$$

which can be identified with the limit $\lim_{\omega \rightarrow \infty} \text{Re} \Sigma_\nu(k, \omega)$, since the dispersive part decays like $1/\omega$ for $\omega \rightarrow \pm\infty$. Eq. (20) looks like a Hartree-Fock (HF) potential. However, $n_\nu(k)$ is the momentum distribution of Eq. (11) containing the depletion effects due to NN correlations and temperature.

III. RESULTS AND DISCUSSION

All the results discussed in this paper have been obtained with the charge-dependent Bonn (CDBONN) potential. Since we want to focus our study on the dependence of the observables on asymmetry, we will consider only one density $\rho = 0.16 \text{ fm}^{-3}$ and one single temperature $T = 5 \text{ MeV}$. This temperature is low enough to allow for conclusions in the limit of zero temperature but high enough to avoid instabilities associated with neutron-proton pairing [17, 18, 19]. As for the proton fractions, we will consider three different cases. The first one corresponds to symmetric nuclear matter $x_p = \rho_p/\rho = 0.50$ and serves as a guide-line for the other cases. The last one has a very low proton fraction $x_p = 0.04$ and corresponds

to the β -stable composition of matter with nucleons, electrons and muons at $T = 5$ MeV and $\rho = 0.16 \text{ fm}^{-3}$. This composition can be computed thanks to the fact that we know the asymmetry dependence of both the neutron and the proton chemical potentials at this density [13]. Finally, we consider an intermediate fraction $x_p = 0.30$ which is useful to identify the effects for low asymmetries of the system.

We start by discussing the momentum dependence of the single-particle spectral functions of neutrons and protons. Depending on how far above or below the considered momentum is from the Fermi momentum of the corresponding particle, we expect a very different behavior for the spectral function. Fig. 1 shows the neutron (left panels) and proton (right panels) single-particle spectral functions for three different momenta ($k = 0, k_F^\nu$ and $2k_F^\nu$, with k_F^ν the Fermi momentum of each nucleon species) at a proton fraction of $x_p = 0.04$. The dotted vertical line corresponds to $\tilde{\omega} \equiv \omega - \mu_\nu = -\mu_\nu$ and indicates the point below which the contribution to the sum rule m_1 in Eq. (18) is negative.

The contributions $A_\nu^<$ and $A_\nu^>$ are given by the dashed and dash-dotted lines, respectively. Since we are considering finite temperature, thermally excited states are always included in the definition of the spectral function Eq. (7). This leads to contributions to $A_\nu^<$ at energies ω larger than μ_ν ($\tilde{\omega}$ larger than zero). Similarly, $A_\nu^>(k, \omega)$ extends to the region below μ_ν . In general there is no longer a clear separation in energy between $A_\nu^>$ and $A_\nu^<$, as it is the case for $T = 0$. Actually, the maxima of $A_\nu^>$ and $A_\nu^<$ can even coincide.

For the case of neutrons at $k = 0$, the peak of the spectral function is provided by $A_n^<$. This is due to the fact that the position of this peak, which can be identified with the quasi-particle energy $\epsilon_{qp}^n(k)$, is well below the chemical potential μ_n , which implies that the value of Fermi-Dirac distribution function $f(\omega)$ at this energy $\omega = \epsilon_{qp}^n(0)$ is very close to one (see Eq. 10). Since the quasi-particle peak is well below μ_n , the thermal effects do not fill up the minimum in the spectral function at $\omega = \mu_n$ and we still observe some kind of separation in energy between $A_n^<$ and $A_n^>$. However, around $\tilde{\omega} = 0$ there is an energy interval, in which both $A_n^<$ and $A_n^>$ are small but different from zero, i.e. an interval in which these functions overlap.

A similar situation is observed for the neutron spectral function at $k = 2k_F^n$. In this case the quasi-particle energy is well above μ_n , which means that $(1 - f(\omega))$ is very close to one at this energy and the peak structure is supplied by $A_n^>$ (see Eq. 10). Therefore we observe a clear separation between the hole part ($A_n^>$) and the particle part ($A_n^<$) of the spectral

function even at finite temperature.

At $k = k_F^n$, the quasi-particle peak is close to the chemical potential. This means that $f(\epsilon_{qp}^n(k))$ is around 0.5 and therefore $A_n^<$ and $A_n^>$ have a peak at $\tilde{\omega} = 0$. The thermal effects fill up the zero that the spectral function has at $\tilde{\omega} = 0$ in the $T = 0$ limit and the overlap region of $A_n^<$ and $A_n^>$ is enhanced.

In the case of very small proton fraction, like $x_p = 0.04$, the Fermi momentum k_F^p is rather small and consequently the quasi-particle energies are close to the Fermi energy μ_p for $k = k_F^p$ as well as $k = 0$ and $k = 2k_F^p$. Therefore, although the relative distances to the Fermi surfaces are the same, there are strong overlaps between $A_p^<$ and $A_p^>$ at all the three momenta. Even in the case of $k = 2k_F^p$, a small peak structure for $A_p^<$ around $\tilde{\omega} = 0$ can be observed.

In Table I (neutrons) and Table II (protons) we report the fraction of the integrated strength of $A_\nu^<$ below and above the corresponding chemical potential, μ_ν . In fact, if we identify $A_\nu^<$ with the $T = 0$ hole spectral function, the integrated strength above the chemical potential would be exactly zero (since $A_\nu^< = 0$ for $\omega > \mu_\nu$). In this sense, the integrated strength above the chemical potential can be considered as a genuine thermal effect. As expected, these thermal effects are important around k_F^ν , where the overlap between $A_\nu^<$ and $A_\nu^>$ is significant. Notice also that for this very neutron-rich system, neutrons are less affected by temperature, while protons (which have a substantially lower Fermi momentum and can thus be considered as a dilute system) are much more influenced by temperature. For instance, protons show a large amount of strength above μ_F^p up to momenta $k/k_F^p \sim 2$. For the sake of completeness, the corresponding occupation numbers for each species are also listed in those tables.

The m_0 sum rule is fulfilled with an accuracy better than .1 % for both neutrons and protons in the complete momentum range. Results for m_1 both for neutrons (upper panels) and protons (lower panels) are reported in Fig. 2 at the three proton fractions $x_p = 0.5, 0.3$ and 0.04. The vertical dotted lines indicate the location of k_F^ν for each case. Since the sum rule is satisfied better than 1 % in all cases, the left- and the right-hand sides of Eq. (18) (solid lines) lie on top of each other and cannot be distinguished. To understand the k -dependence of the different contributions to m_1 , it is useful to keep in mind the proton and neutron chemical potentials at the different concentrations given in Table III.

The lower dash-dotted line shows the m_1 contribution from $A_\nu^<$. For momenta below k_F^ν , the

contribution of $A_\nu^<$ to m_1 is dominated by the quasi-particle peak, which lies below μ_ν . As the momentum increases, the peak appears closer to $\omega = 0$ and its weight in the integral is diminished, which makes the integral smaller in magnitude and thus the contribution to the sum rule becomes an increasing function of k . When we get closer to the Fermi momentum k_F^ν , the peak moves to $\omega \sim \mu_\nu$ and the position of the chemical potential becomes crucial. As far as the chemical potential is negative, the contribution of $A_\nu^<$ will be negative because the Fermi-Dirac function falls off close to μ_ν and the integral will only catch a little positive zone if the chemical potential is close to $\omega \sim 0$. If, on the other hand, the chemical potential is positive, the integrand is not zero for $\omega > 0$ and there can be a non-negligible positive contribution to the integral when the quasi-particle peak lies between $\omega = 0$ and $\omega = \mu_\nu$. This cancellations between positive and negative contributions give rise to the structures observed for the $A_\nu^<$ sum rule close to the Fermi momentum. In the upper-right panel of Fig. 2, for instance, the asymmetry is so extreme that the neutron chemical potential $\mu_n > 0$ and the positive contribution to the integral is important enough to pull the sum rule to positive values for $k \sim k_F^n$. Finally, for high momenta, $A_\nu^<$ is strongly suppressed (the quasi-particle peak lies in the region $\omega > \mu_\nu$, which is suppressed by the Fermi-Dirac factor) and the contribution to the sum rule goes to zero.

The upper dash-dotted line displays the contribution from $A_\nu^>$. Due to the short-range correlations, there is always a high energy tail that gives rise to a positive contribution which, for momenta well below k_F^ν , is nearly constant. In the case of the neutron sum rules, this constant decreases slightly with decreasing proton fraction, although the chemical potential μ_n increases significantly. Such a small decrease is in accordance with the fact that the neutron-proton correlations are dominant compared to the neutron-neutron ones. These high energy tails (which are mainly caused by short-range correlations) are, nevertheless, not much affected by asymmetry. When the momenta becomes larger than k_F^n , m_1 increases monotonously following the location of the quasi-particle peak, which moves to higher energies when the momentum grows.

Also for protons one observes a contribution from $A_\nu^>$ to the sum rule which is almost a constant for momenta smaller than the Fermi momentum. In this case, however, this constant is almost independent of the proton fraction, although the chemical potential μ_p gets significantly more attractive with decreasing x_p . At small values of x_p , the neutron abundance is large and this leads to strong correlation effects in the proton spectral function. In

particular at $x_p = 0.04$, which is the case corresponding to the lowest Fermi momenta, the total contribution of $A_p^>$ is the result of a balance between positive and negative contributions. These can result in a minimum for $k \sim k_F^p$. In that particular proton fraction (see Fig. 1), the quasi-particle peaks lies always below $\omega = 0$ and thus it gives rise to a negative contribution. The relative width of this peak and the detailed structure of the high-energy tail can push the integral to negative values (almost -15 MeV at $k \sim 160$ MeV). At larger momenta, however, the contribution to m_1 of $A_p^>$ starts to grow steadily. Notice that for this asymmetry we can not say that this is due to the movement of the quasi-particle peak, because even at $k = 2k_F^p$ the quasi-particle peak lies in the region where the contribution to the integral is negative. The growth in k is, to a large extent, a consequence of the contributions to the spectral functions at large energies.

In addition to these contributions, we have plotted in the dotted lines of Fig. 2 the Hartree-Fock approximation of m_1 at the same temperature, density and proton fraction. This approximation turns out to give a very good estimate of m_1 , an interesting result which has been already observed in the previous analysis for symmetric nuclear matter [14, 15]. This result allows for a quantitative estimate of the amount of correlations produced by a given NN potential without performing sophisticated many-body calculations.

Furthermore, it is worth noticing that the value of m_1 (specially for large k) is not so much affected by the asymmetry of the system. It is very similar for both neutrons and protons in all the three cases considered. This is a somehow surprising result, since we have seen that the separate contributions of $A_\nu^<$ and $A_\nu^>$ can indeed be very different. Of course, at high momentum the kinetic energy is dominant and this could explain this fact in part.

However, even at zero momentum the effect of the asymmetry is rather moderate, as we can see from Fig. 3, where Σ_ν^∞ , Σ_ν^{HF} (the Hartree-Fock approximation to the self-energy) and the quasi-particle energy ϵ_{qp}^ν at zero momentum are plotted as a function of the proton fraction. Let us consider the highly asymmetric case of $x_p = 0.04$. The isospin splitting of the Hartree-Fock energies ($k = 0$) at this asymmetry is around 10 MeV, which is rather small compared to the isospin splitting of the quasi-particle energies, which is around 55 MeV. This means that the stronger binding of the proton states as compared to the neutron states is to a small extent due to the attractive neutron-proton interaction in the bare NN potential, the Born term of the T -matrix. The obtained attraction is then mainly due to the terms in the T -matrix coming from the second and higher orders in the NN potential. In

other words, it is an effect of strong correlations in the isospin equal to zero NN channels (like the ${}^3S_1 - {}^3D_1$ partial wave), which lead to the deeply bound quasi-particle energy for protons at this large neutron abundance. These correlation effects, however, are also important for a redistribution of single-particle strength for protons with $k = 0$ to energies well above the Fermi energy, leading to a low occupation probability (see Table II) and a large positive contribution to the sum rule m_1 . This positive contribution yields to a value of Σ^∞ , which is even above the HF energy. The effects of correlations for the corresponding energies of the neutron state are much weaker, indicating that correlation effects in the isospin one NN channels are weaker.

So we observe a compensation of correlation effects in the energy weighted sum rule. On the one hand, correlations lead to a more attractive quasi-particle energy (compared to the HF result) and therefore to a more attractive contribution of the quasi-particle pole to the sum rule. On the other hand, correlations shift single-particle strength to high energies, which yield a repulsive contribution to m_1 . This compensation of correlation effects explains that the value of the energy integrated sum rule for the correlated system can be very similar to the HF result.

This compensation of correlation effects can also be observed in Fig. 4, which shows the exhaustion of the sum rules m_0 (left panel) and m_1 (right panel) at a momentum $k = 400$ MeV for both neutrons (solid lines) and protons (dashed lines), at the three different proton fractions. The first thing to be observed is that the asymmetry affects mainly the amount of strength exhausted at intermediate energy regions, while there is not a noticeable difference at high energies. For $x_p = 0.5$, i.e. the symmetric case, the quasi-particle energies for neutrons and protons at $k = 400$ MeV, are the same: $\epsilon_{qp} = 36.45$ MeV. When the proton fraction decreases, the quasi-particle energies split. Neutrons become more repulsive ($\epsilon_{qp}^n = 41.88$ MeV), in contrast with protons ($\epsilon_{qp}^p = 30.86$ MeV). This behavior is confirmed at $x_p = 0.04$, resulting in $\epsilon_{qp}^n = 49.35$ MeV and $\epsilon_{qp}^p = 25.22$ MeV. In the case of m_0 , since the quasi-particle peak of protons is located at lower energies, the amount of strength exhausted at lower energies is larger for protons than for neutrons. This trend changes at intermediate energies and both strengths merge together for large energies, which is an indication of the fact that, for these momenta and asymmetries, the quasi-particle contributions to the sum rule of neutrons are larger than those of protons, thus explaining why the increase in m_0 at the quasi-particle pole is larger in the former case. The same type of analysis is valid for

m_1 . Notice that, as previously mentioned, the final value of m_1 is not much affected by the asymmetry and the behavior at high energies is the same for all the cases.

To summarize, we have analyzed the behavior of the energy weighted sum rules of single-particle spectral functions of hot asymmetric nuclear matter. The sum rules are very well fulfilled, because the T -matrix approximation itself respects the analytical properties of both the self-energy and the Green's function, in which the sum rules are based. Nevertheless, they are a good test of the numerical consistency of the calculation, which may be helpful e.g. in deciding the best distribution of the energy mesh points when one has to work with spectral functions.

Employing realistic NN interactions, an important source of correlations come from the strong components in the neutron-proton interaction. As a consequence, one observes in neutron-rich matter a larger depletion of the occupation probabilities for protons with momenta below the Fermi-momentum than for neutrons. One also finds that these correlations, at this asymmetry, lead to much more attractive quasi-particle energies for protons than those obtained in the HF approximation. The same effect is observed in neutrons, although it is considerably weaker. This shift of the quasi-particle energies tends to lead to more attractive energy weighted sum rules m_1 . This effect is compensated by the fact that the very same correlations are also responsible for a shift of single-particle strength to high positive energies. As a consequence, the energy integrated sum rules and the isospin splitting in asymmetric nuclear matter for the correlated system yield results which are very close to the HF ones. In both cases, however, there is not a strong dependence on asymmetry due to the lack of tensor correlations. This relocation of single-particle strength, on the other hand, can be nicely observed in the convergence of the energy weighted sum rule, which is therefore an indicator of correlation effects.

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k/k_F^n	below μ_n [%]	above μ_n [%]	$n(k)$
0.0	99.98	0.02	0.971
0.5	99.96	0.04	0.960
1.0	55.58	44.42	0.464
1.5	98.48	1.52	0.006
2.0	99.76	0.24	0.001

TABLE I: Strength distribution of $A_n^<$ at $\rho = 0.16 \text{ fm}^{-3}$, $T = 5 \text{ MeV}$ and proton fraction $x_p = 0.04$. The numbers give the fraction of the integrated strength above and below the neutron chemical potential μ_n . The last column reports the occupation of the respective neutron momentum state.

k/k_F^p	below μ_p [%]	above μ_p [%]	$n(k)$
0.0	97.62	2.38	0.605
0.5	94.88	5.12	0.531
1.0	21.62	78.38	0.280
1.5	24.26	75.74	0.056
2.0	64.10	35.90	0.009

TABLE II: Strength distribution of $A_p^<$ at $\rho = 0.16 \text{ fm}^{-3}$, $T = 5 \text{ MeV}$ and proton fraction $x_p = 0.04$. The numbers give the fraction of the integrated strength above and below the proton chemical potential μ_p . The last column reports the occupation of the respective proton momentum state.

x_p	μ_n [MeV]	μ_p [MeV]
0.50	-23.46	-23.46
0.30	-2.78	-48.33
0.04	19.21	-91.45

TABLE III: Neutron and proton chemical potentials at $\rho = 0.16 \text{ fm}^{-3}$, $T = 5 \text{ MeV}$ and different proton fractions.

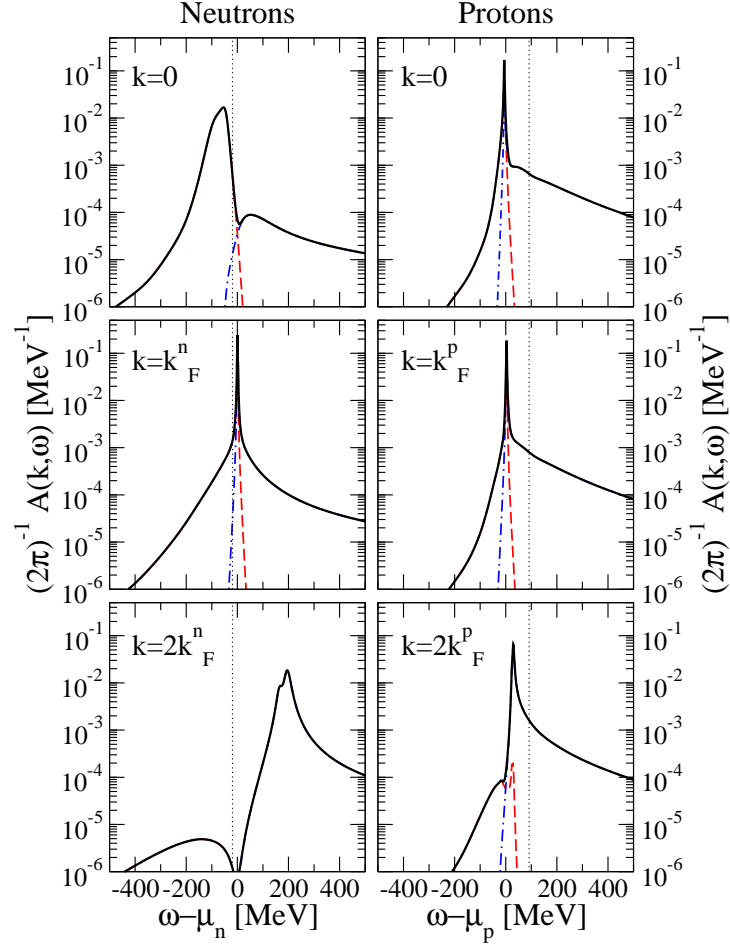


FIG. 1: (Color online) Neutron (left panels) and proton (right panels) single-particle spectral functions at $\rho = 0.16 \text{ fm}^{-3}$, $T = 5 \text{ MeV}$ and proton fraction $x_p = 0.04$ for three different momenta. Both $A_\nu^<$ (dashed lines) and $A_\nu^>$ (dot-dashed lines) are displayed together with the total spectral function A_ν (solid lines).

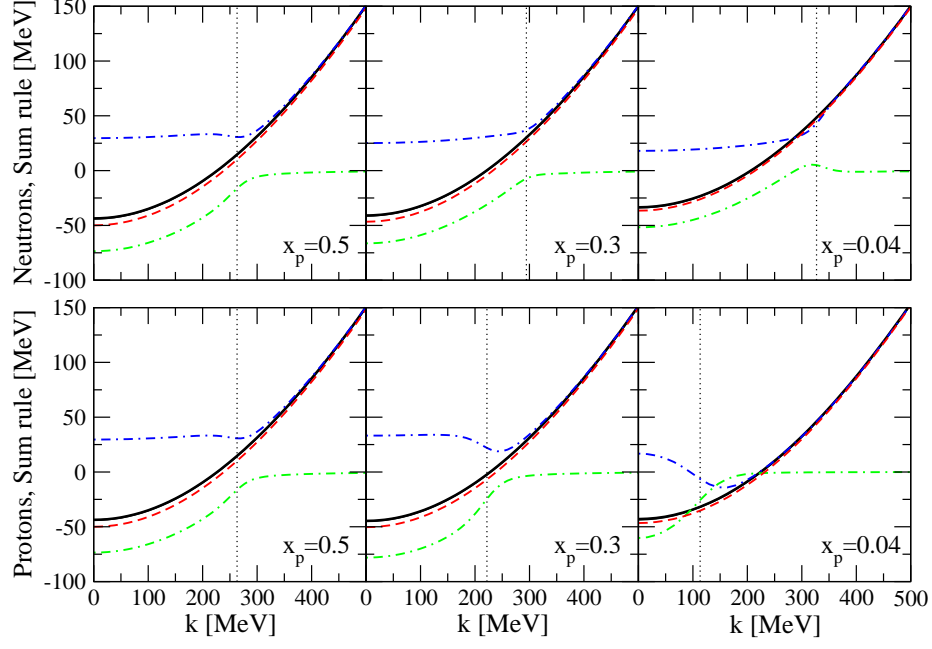


FIG. 2: (Color online) Energy weighted sum rule m_1 (solid lines) for neutrons (panels above) and protons (panels below) at $\rho = 0.16 \text{ fm}^{-3}$, $T = 5 \text{ MeV}$ and several proton fractions. Both the right- and the left-hand sides of Eq. (18) are displayed, but the sum rule is so well fulfilled that they cannot be distinguished. The contributions to m_1 that come from $A_\nu^<$ and $A_\nu^>$ are indicated by the upper and the lower dash-dotted lines, respectively. The result for the energy weighted sum rule obtained from the Hartree-Fock approximation is represented by the dashed lines.

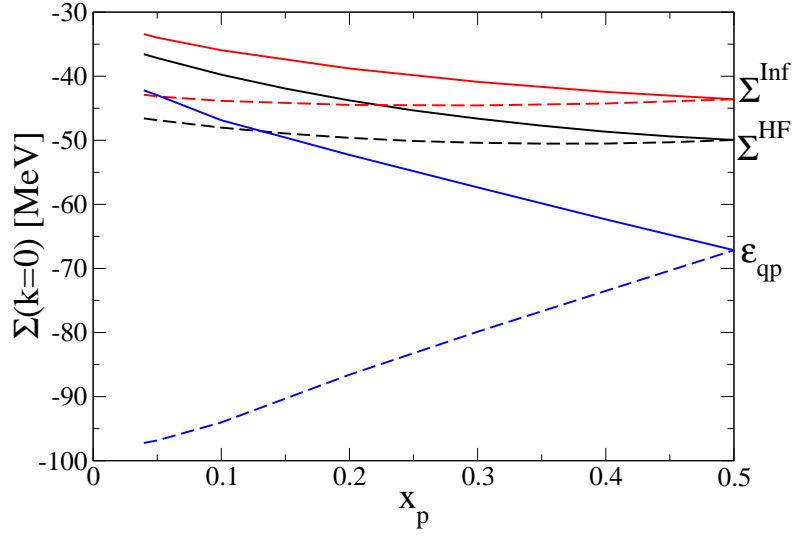


FIG. 3: (Color online) Σ_ν^∞ , Σ_ν^{HF} and ϵ_{qp}^ν at zero momentum as a function of the proton fraction at $\rho = 0.16 \text{ fm}^{-3}$ and $T = 5 \text{ MeV}$ for neutrons (solid lines) and protons (dashed lines).

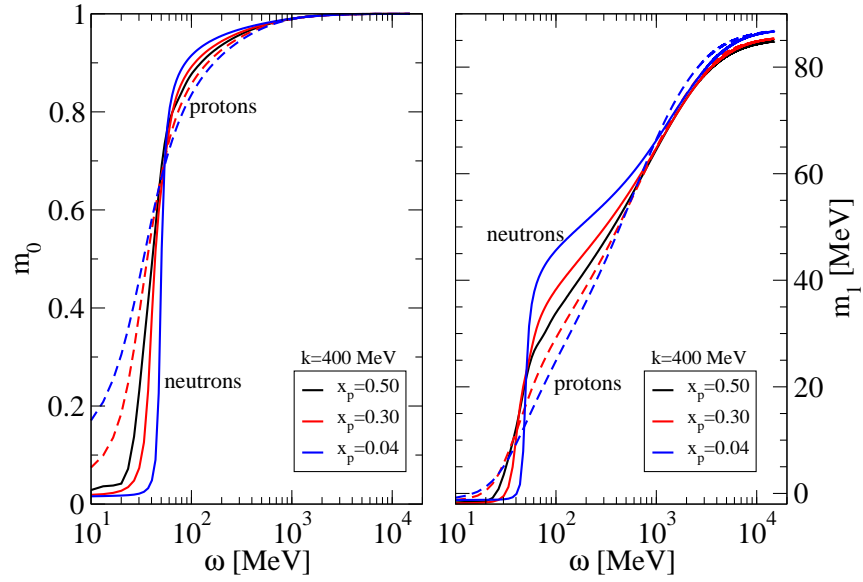


FIG. 4: (Color online) Saturation of the sum rules m_0 (right panel) and m_1 (left panel) for neutrons (solid lines) and protons (dashed lines) at three different proton fractions. The momentum is $k=400 \text{ MeV}$ and the density and temperature, the same same as in the previous figures.